LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120619\
Data File : BM024017.D
Acq On : 06 Dec 2019 15:42
Operator : JU
Sample : K6007-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BL5

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.034	23	25	38	rVB	56943	109546	1.79%	0.120%
2	3.534	107	110	117	rVB	64749	86072	1.40%	0.094%
3	3.651	125	130	139	rVB2	64369	120636	1.97%	0.132%
4	3.734	139	144	154	rVB	116126	177552	2.90%	0.194%
5	4.646	294	299	307	rVV2	58353	155844	2.54%	0.171%
6	5.151	380	385	397	rVB	167432	314821	5.14%	0.345%
7	5.522	441	448	453	rBV2	95732	161101	2.63%	0.176%
8	6.175	554	559	566	rVB2	99041	155922	2.54%	0.171%
9	6.475	605	610	617	rVB4	53392	108858	1.78%	0.119%
10	6.669	632	643	656	rVV	861034	1678178	27.39%	1.838%
11	6.828	665	670	680	rVV	705516	1181838	19.29%	1.294%
12	6.916	680	685	690	rVV	122463	202289	3.30%	0.222%
13	7.016	696	702	714	rVB	965162	1545172	25.22%	1.692%
14	7.122	714	720	729	rBV2	98097	206349	3.37%	0.226%
15	7.381	756	764	772	rVV2	85578	163584	2.67%	0.179%
16	7.475	774	780	790	rVV	1534193	2452148	40.02%	2.686%
17	7.698	812	818	823	rVV2	61646	104945	1.71%	0.115%
18	8.198	897	903	914	rBV	895551	1535163	25.05%	1.681%
19	8.304	914	921	928	rVV3	51874	133969	2.19%	0.147%
20	8.628	970	976	985	rVV	819001	1404269	22.92%	1.538%
21	8.710	986	990	998	rVV	102622	164687	2.69%	0.180%
22	9.345	1091	1098	1110	rBV	661916	1177769	19.22%	1.290%
23	9.669	1148	1153	1163	rVB	289466	498331	8.13%	0.546%
24	9.880	1181	1189	1202	rVV	1037813	1872240	30.55%	2.051%
25	10.245	1241	1251	1258	rBV	2158641	3612191	58.95%	3.957%
26	10.298	1258	1260	1269	rVB	71085	118496	1.93%	0.130%
27	10.398	1269	1277	1291	rBV	568721	1215307	19.83%	1.331%
28	11.298	1423	1430	1435	rBV	83086	141397	2.31%	0.155%
29	11.727	1491	1503	1511	rBV	3860420	6127759	100.00%	6.712%
30	11.922	1529	1536	1545	rVB	199989	344441	5.62%	0.377%
31	12.145	1568	1574	1581	rBV	107322	187197	3.05%	0.205%
32	12.974	1704	1715	1717	rVV2	205276	349782	5.71%	0.383%
33	12.998	1717	1719	1726	rVV	175904	276562	4.51%	0.303%
34	13.298	1764	1770	1771	rBV2	79342	115517	1.89%	0.127%

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Misc :
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Instrument : BNA_M
ClientSampleId : C0BL5

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Title : SVOA CALIBRATION

35	13.457	1792	1797	1802	rVV	142803	259482	4.23%	0.284%
36	13.510	1802	1806	1809	rVV	119918	184612	3.01%	0.202%
37	13.563	1809	1815	1819	rVV	1938220	2837161	46.30%	3.108%
38	13.604	1819	1822	1832	rVV	723972	1122845	18.32%	1.230%
39	13.692	1832	1837	1842	rVV2	85338	150057	2.45%	0.164%
40	13.827	1853	1860	1871	rBV	2313029	3439081	56.12%	3.767%
41	14.063	1892	1900	1905	rBV	188323	249722	4.08%	0.274%
42	14.139	1905	1913	1920	rBV2	3096318	4831728	78.85%	5.292%
43	14.233	1925	1929	1933	rVV3	76986	142321	2.32%	0.156%
44	14.374	1946	1953	1960	rBV3	391284	804519	13.13%	0.881%
45	14.545	1972	1982	1986	rBV3	136349	325260	5.31%	0.356%
46	14.586	1986	1989	1992	rVV4	94103	141226	2.30%	0.155%
47	14.627	1992	1996	2002	rVB	90426	135514	2.21%	0.148%
48	14.762	2014	2019	2025	rBV2	76278	140382	2.29%	0.154%
49	14.821	2025	2029	2034	rVB	98305	129939	2.12%	0.142%
50	14.939	2043	2049	2052	rBV2	59537	96149	1.57%	0.105%
51	15.027	2059	2064	2070	rVB	216388	271871	4.44%	0.298%
52	15.139	2077	2083	2089	rVV	3058131	4279364	69.84%	4.687%
53	15.192	2089	2092	2098	rVV4	57593	109593	1.79%	0.120%
54	15.280	2103	2107	2120	rVV	636869	1000267	16.32%	1.096%
55	15.433	2127	2133	2137	rVB5	44940	97588	1.59%	0.107%
56	15.492	2137	2143	2154	rVB3	135548	267318	4.36%	0.293%
57	15.639	2162	2168	2177	rVB5	62147	162808	2.66%	0.178%
58	15.727	2177	2183	2188	rVB4	60844	108676	1.77%	0.119%
59	15.886	2206	2210	2223	rVB	284921	461590	7.53%	0.506%
60	16.251	2269	2272	2277	rVB	68098	87460	1.43%	0.096%
61	16.557	2319	2324	2331	rVB4	108064	191436	3.12%	0.210%
62	16.633	2331	2337	2342	rBV3	78438	175284	2.86%	0.192%
63	16.680	2342	2345	2349	rBV	248067	301083	4.91%	0.330%
64	16.892	2375	2381	2385	rBV	3651326	5049658	82.41%	5.531%
65	16.933	2385	2388	2392	rVV	548131	711839	11.62%	0.780%
66	16.992	2392	2398	2411	rVB	3065704	4322524	70.54%	4.735%
67	17.092	2411	2415	2420	rBV5	48515	84773	1.38%	0.093%
68	17.215	2432	2436	2441	rVB2	92481	133453	2.18%	0.146%
69	17.421	2467	2471	2475	rBV	221769	270828	4.42%	0.297%
70	17.768	2524	2530	2534	rBV2	212468	371793	6.07%	0.407%
71	17.815	2534	2538	2549	rVV	971957	1373411	22.41%	1.504%

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Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Title : SVOA CALIBRATION

72	17.956	2557	2562	2566	rBV2	189878	288137	4.70%	0.316%
73	17.998	2566	2569	2574	rVB	143278	212942	3.48%	0.233%
74	18.115	2581	2589	2594	rBV	240285	406367	6.63%	0.445%
75	18.274	2613	2616	2621	rVB	142469	204622	3.34%	0.224%
76	18.580	2664	2668	2671	rBV	80556	107565	1.76%	0.118%
77	18.698	2684	2688	2693	rVV	180368	264948	4.32%	0.290%
78	18.756	2693	2698	2702	rVV	272774	364873	5.95%	0.400%
79	18.792	2702	2704	2708	rVB	72052	84146	1.37%	0.092%
80	18.839	2708	2712	2721	rBV3	67340	158524	2.59%	0.174%
81	18.962	2729	2733	2742	rBV2	349832	673102	10.98%	0.737%
82	19.033	2742	2745	2749	rVB	71700	91192	1.49%	0.100%
83	19.109	2754	2758	2765	rBV	269830	423836	6.92%	0.464%
84	19.297	2785	2790	2796	rBV	3328024	4723816	77.09%	5.174%
85	19.409	2807	2809	2815	rVB3	81961	131439	2.14%	0.144%
86	19.662	2848	2852	2854	rBV3	67983	97146	1.59%	0.106%
87	19.833	2878	2881	2885	rVB	84836	110649	1.81%	0.121%
88	19.880	2886	2889	2895	rVB	202155	225734	3.68%	0.247%
89	20.256	2949	2953	2959	rVB	336981	378810	6.18%	0.415%
90	20.380	2971	2974	2978	rBV	200666	213163	3.48%	0.233%
91	20.844	3049	3053	3057	rBV	567415	707229	11.54%	0.775%
92	21.103	3091	3097	3101	rBV	3799404	5144574	83.96%	5.635%
93	21.280	3124	3127	3131	rBV	292957	362336	5.91%	0.397%
94	21.703	3196	3199	3204	rBV2	332820	376549	6.14%	0.412%
95	22.144	3271	3274	3281	rVB	387421	520604	8.50%	0.570%
96	22.627	3353	3356	3362	rVB2	506667	691386	11.28%	0.757%
97	23.109	3432	3438	3443	rBV	2606309	4460306	72.79%	4.885%
98	23.162	3444	3447	3453	rVB2	494750	775769	12.66%	0.850%
99	23.244	3455	3461	3470	rVB	2928177	5116435	83.50%	5.604%
100	23.774	3547	3551	3558	rVB	547652	958729	15.65%	1.050%

Sum of corrected areas: 91297475

LSC Report - Integrated Chromatogram

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Data File : BM024017.D
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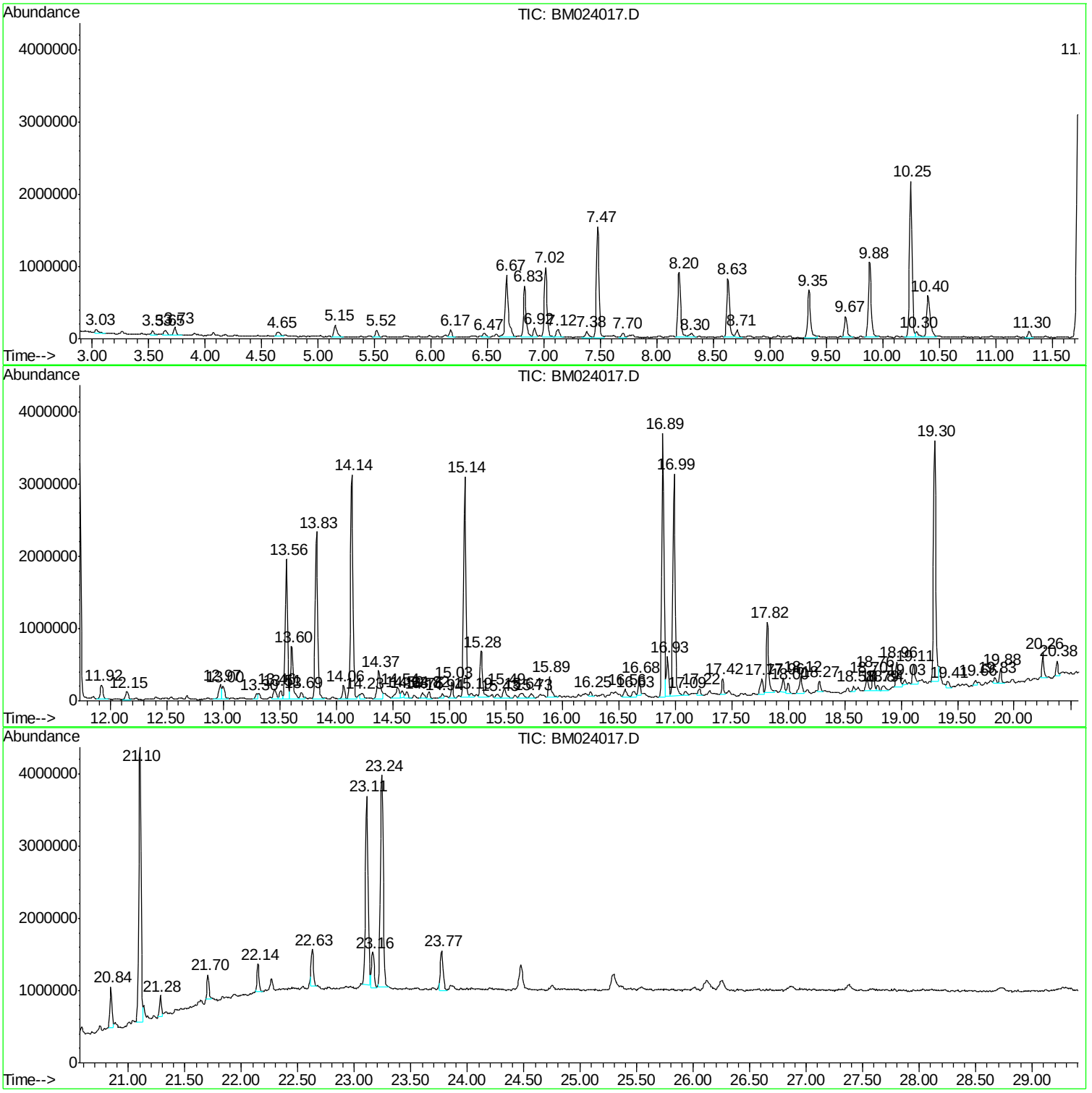
Sample : K6007-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BL5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120619\
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Quant Title : SVOA CALIBRATION

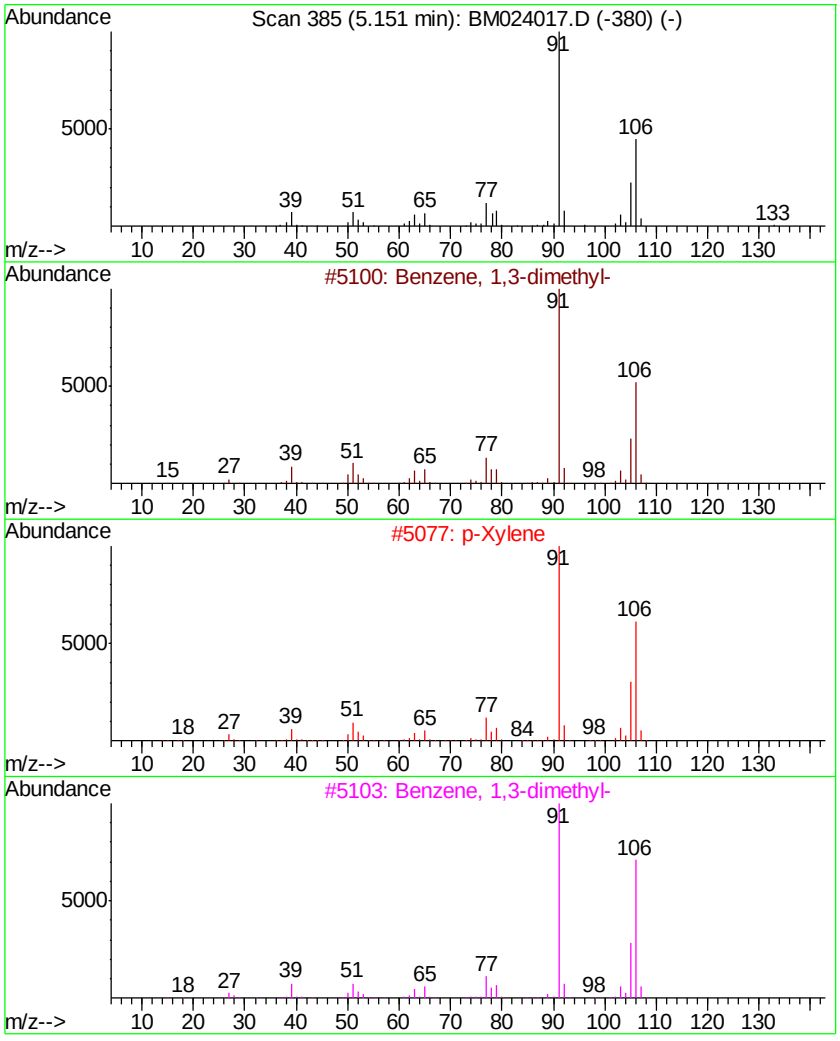
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

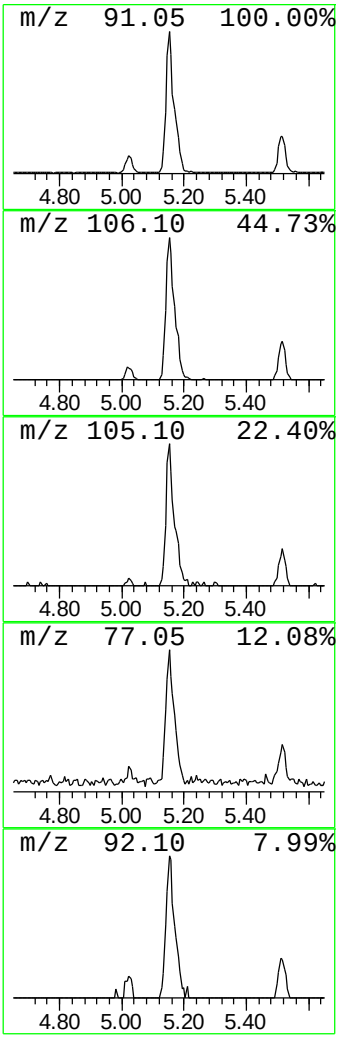
****

Peak Number 1 Benzene, 1,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	2.57 ng/ul	314821	1,4-Dichlorobenzene-d4	7.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2	p-Xylene	106	C8H10	000106-42-3	95
3	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95
4	p-Xylene	106	C8H10	000106-42-3	95
5	p-Xylene	106	C8H10	000106-42-3	95





Library Search Compound Report

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ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BL5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

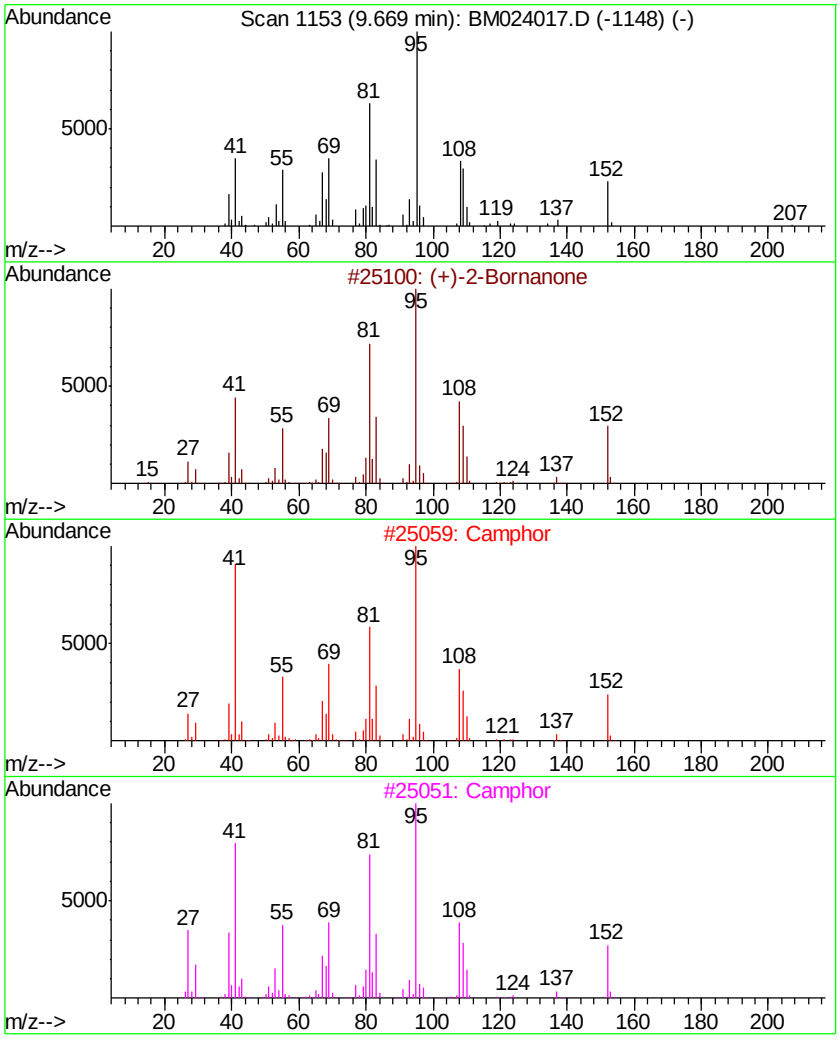
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

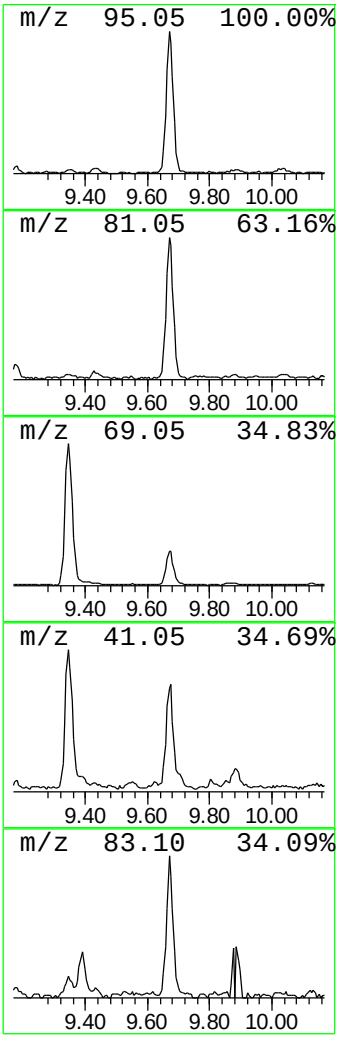
****

Peak Number 2 (+)-2-Bornanone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.67	2.76 ng/ul	498331	Naphthalene-d8	10.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+)-2-Bornanone	152	C10H16O	000464-49-3	97
2	Camphor	152	C10H16O	000076-22-2	97
3	Camphor	152	C10H16O	000076-22-2	96
4	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	96
5	(+)-2-Bornanone	152	C10H16O	000464-49-3	96





Library Search Compound Report

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Quant Title : SVOA CALIBRATION

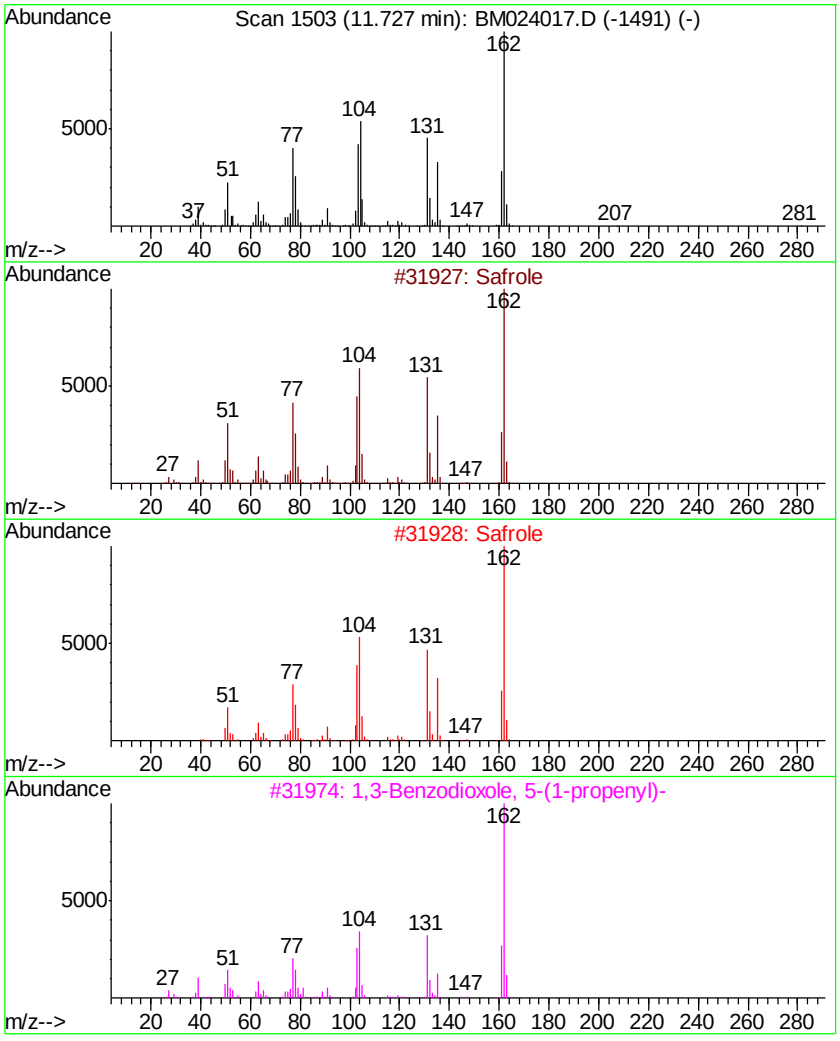
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

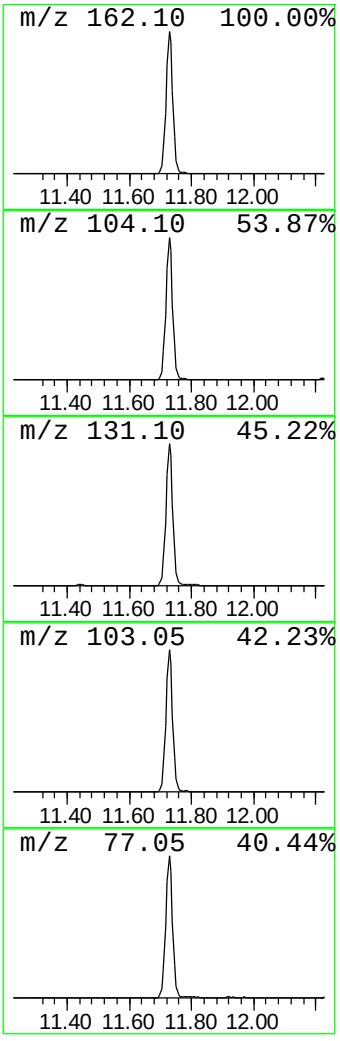
****

Peak Number 3 Safrole Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.73	33.93 ng/ul	6127760	Naphthalene-d8	10.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Safrole	162	C10H10O2	000094-59-7	98
2	Safrole	162	C10H10O2	000094-59-7	98
3	1,3-Benzodioxole, 5-(1-propenyl)-	162	C10H10O2	000120-58-1	97
4	Safrole	162	C10H10O2	000094-59-7	96
5	1,3-Benzodioxole, 5-(1-propenyl)-	162	C10H10O2	000120-58-1	96





Library Search Compound Report

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Instrument : BNA_M
ClientSampleId : C0BL5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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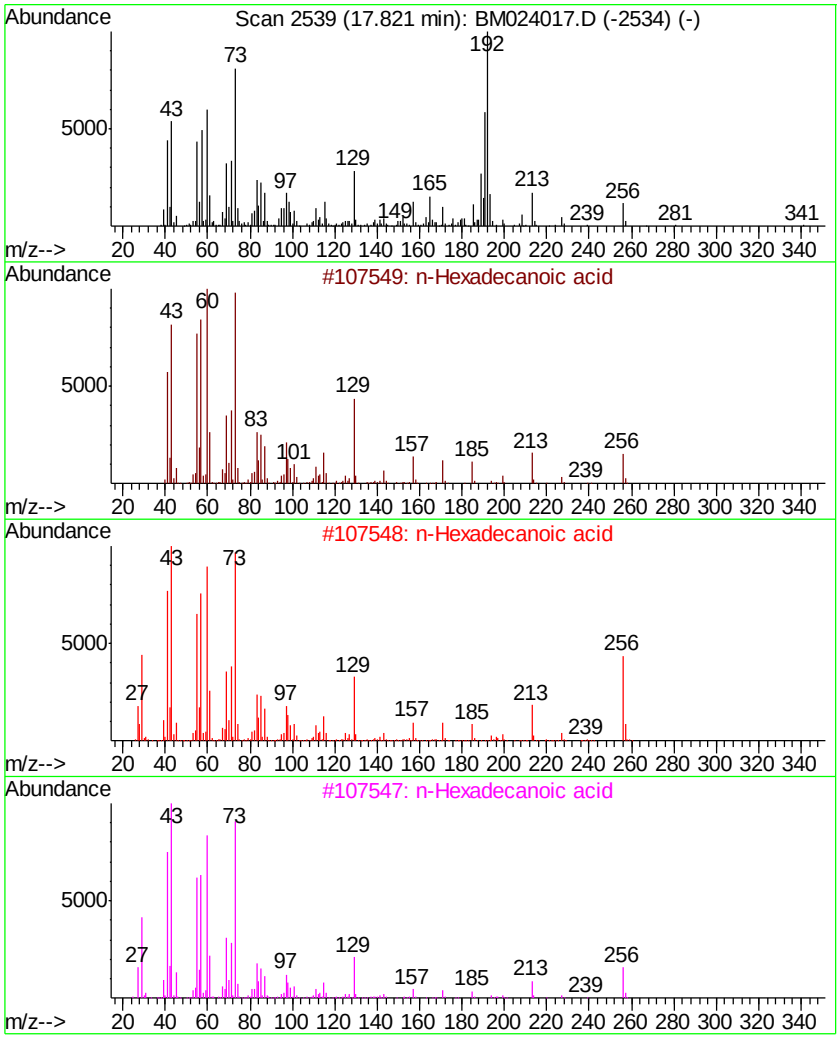
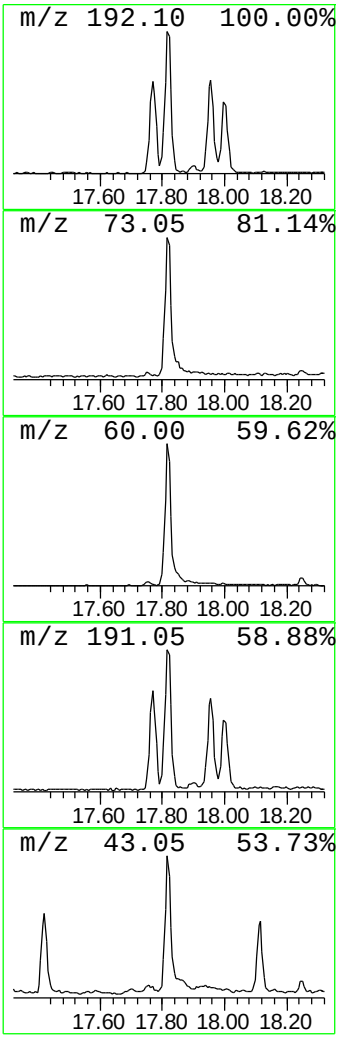
Peak Number 4 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.82	5.44 ng/ul	1373410	Phenanthrene-d10	16.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	89





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Sample : K6007-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BL5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

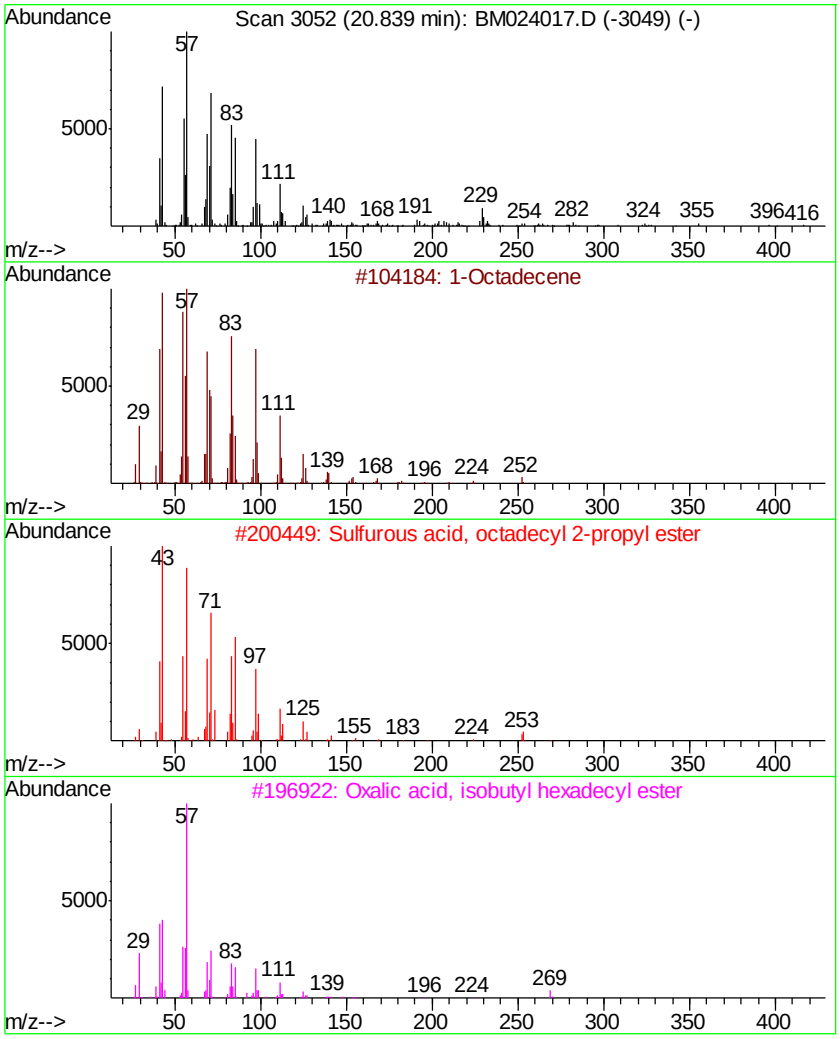
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

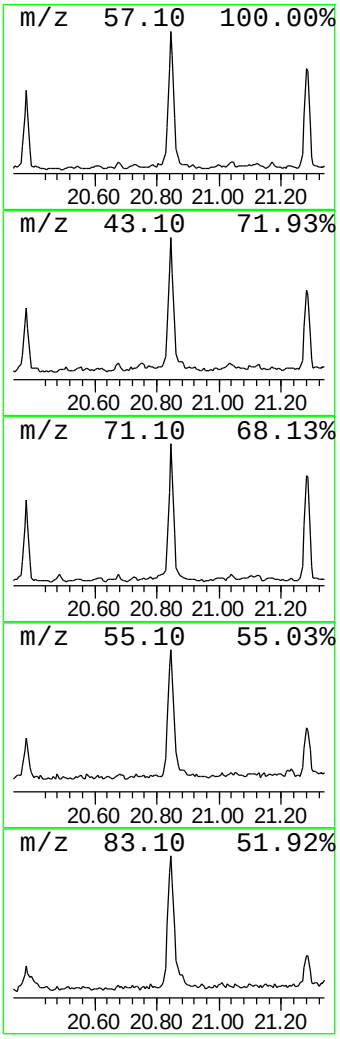
****

Peak Number 5 (DEL) Alkane: Cyclic20.84 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.84	2.75 ng/ul	707229	Chrysene-d12	21.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Octadecene	252	C18H36	000112-88-9	95
2	Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	91
3	Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	91
4	Oxalic acid, isobutyl heptadecyl...	384	C23H44O4	1000309-38-2	91
5	Ethanol, 2-(hexadecyloxy)-	286	C18H38O2	002136-71-2	87





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120619\
Data File : BM024017.D
Acq On : 06 Dec 2019 15:42
Operator : JU
Sample : K6007-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BL5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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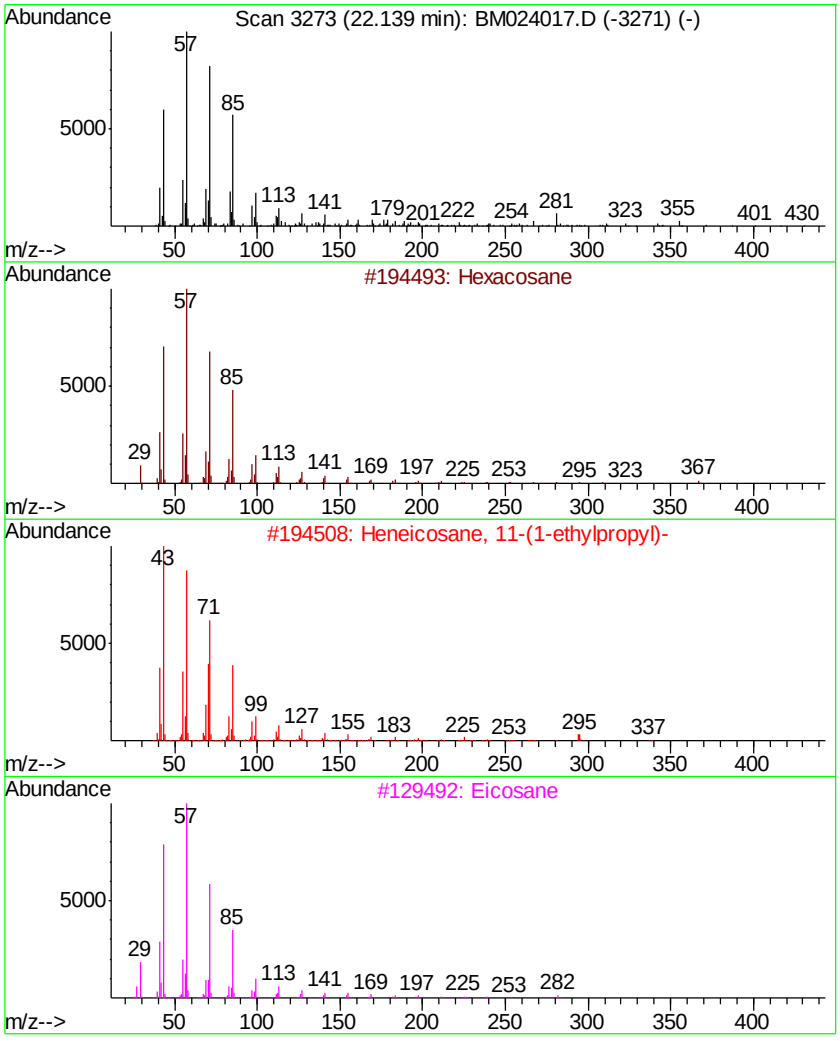
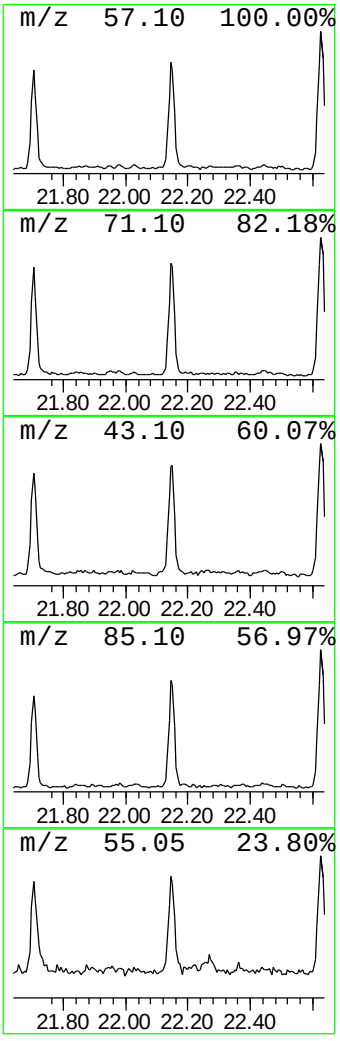
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.14	2.02 ng/ul	520604	Chrysene-d12	21.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	91
2	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
3	Eicosane	282	C20H42	000112-95-8	90
4	Tetracosane	338	C24H50	000646-31-1	90
5	Tetradecane	198	C14H30	000629-59-4	87





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Acq On : 06 Dec 2019 15:42
Operator : JU
Sample : K6007-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BL5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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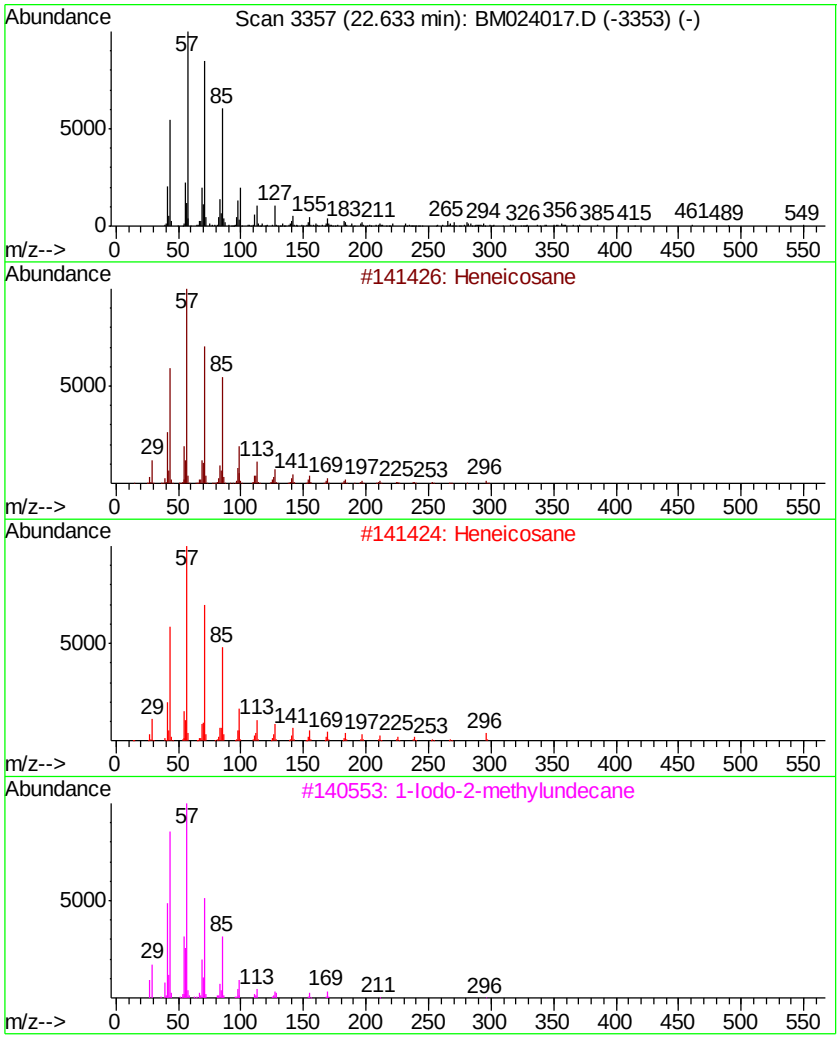
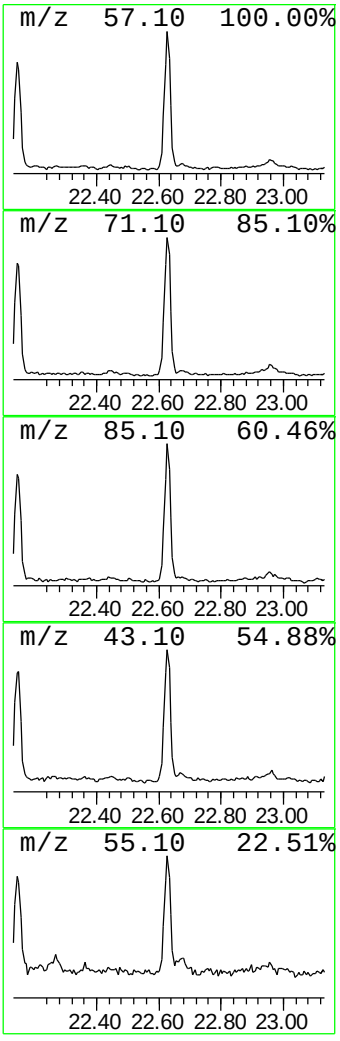
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.63	2.70 ng/ul	691386	Perylene-d12	23.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	94
2	Heneicosane	296	C21H44	000629-94-7	93
3	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	91
4	Hexacosane	366	C26H54	000630-01-3	91
5	Triacontane	422	C30H62	000638-68-6	91





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120619\
Data File : BM024017.D
Acq On : 06 Dec 2019 15:42
Operator : JU
Sample : K6007-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BL5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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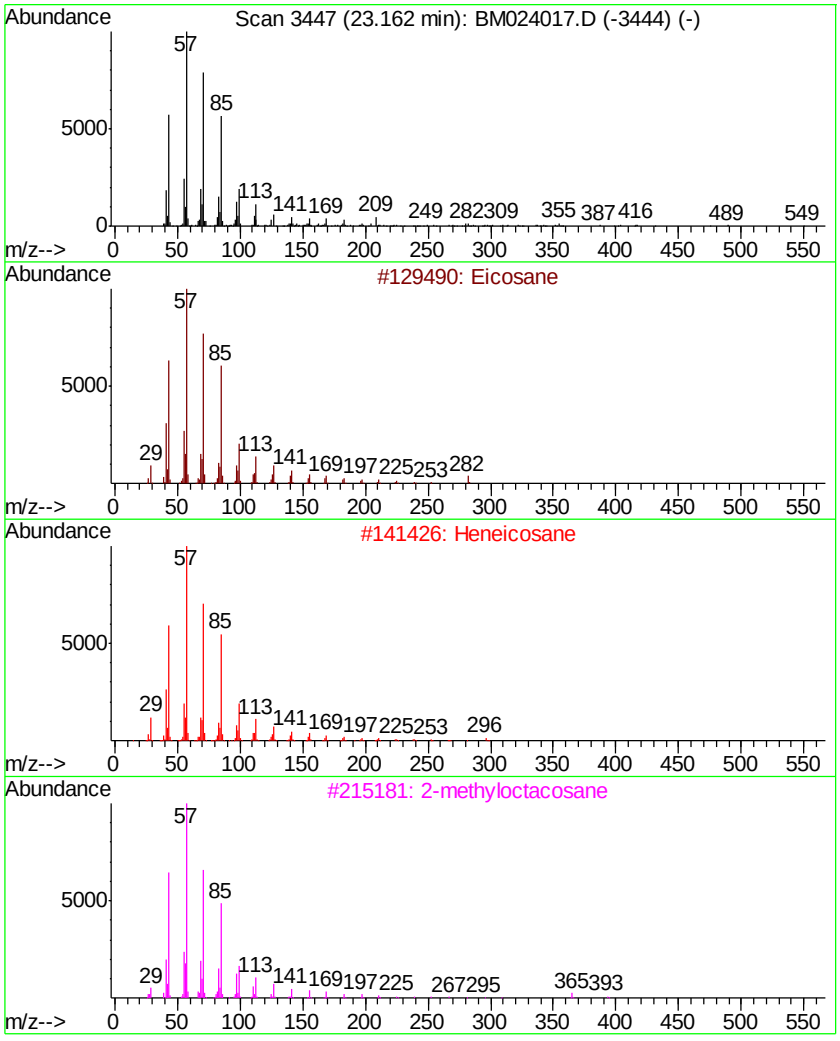
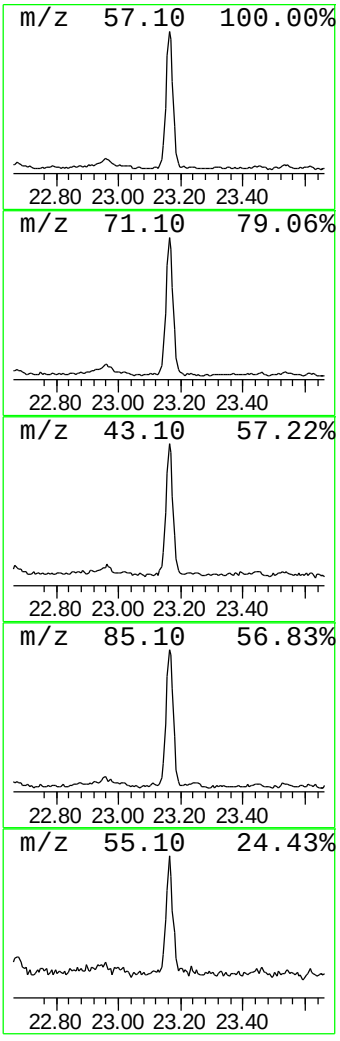
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.16	3.03 ng/ul	775769	Perylene-d12	23.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	97
2	Heneicosane	296	C21H44	000629-94-7	91
3	2-methyloctacosane	408	C29H60	1000376-72-8	91
4	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	90
5	Hexacosane	366	C26H54	000630-01-3	90





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120619\
Data File : BM024017.D
Acq On : 06 Dec 2019 15:42
Operator : JU
Sample : K6007-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BL5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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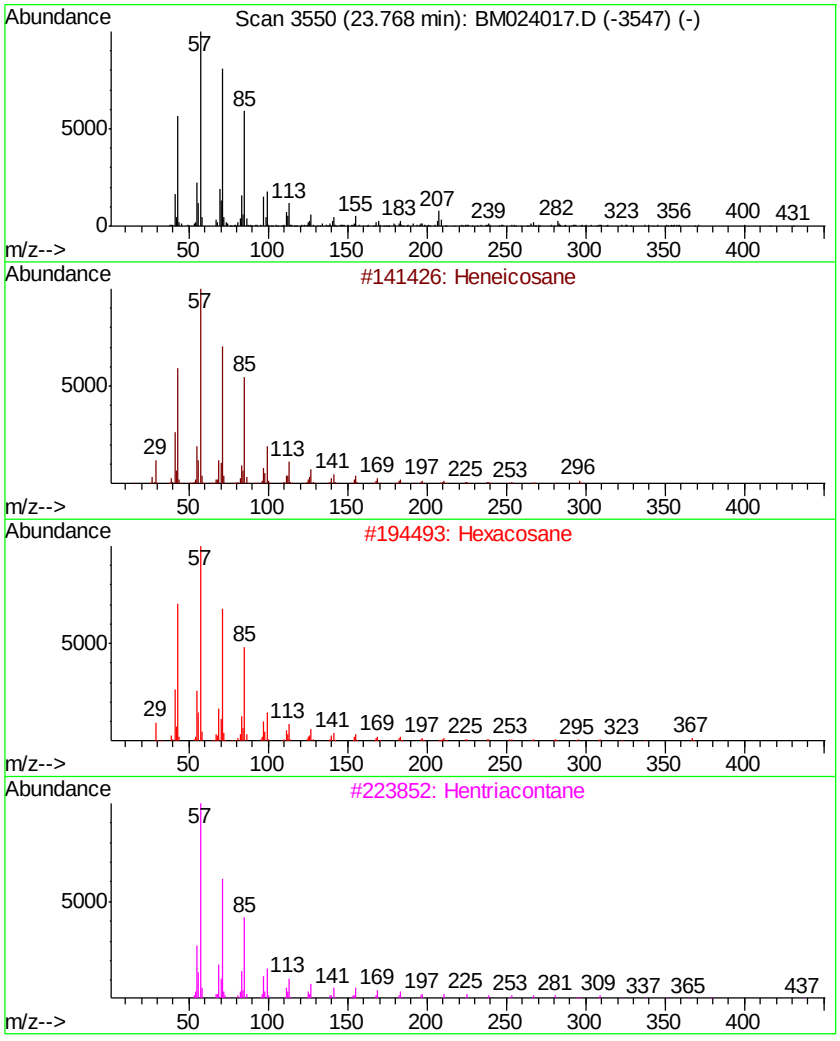
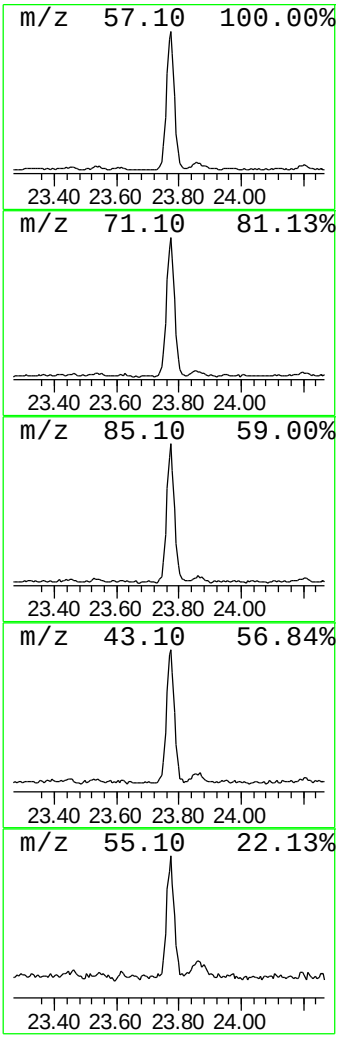
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.77	3.75 ng/ul	958729	Perylene-d12	23.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Hexacosane	366	C26H54	000630-01-3	91
3	Hentriacontane	437	C31H64	000630-04-6	91
4	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90
5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM120619\
Data File : BM024017.D
Acq On : 06 Dec 2019 15:42
Operator : JU
Sample : K6007-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0BL5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, 1,3-dime...	5.15	2.6	ng/ul	314821	1	7.47	2452150	20.0
(+)-2-Bornanone	9.67	2.8	ng/ul	498331	2	10.25	3612190	20.0
Safrole	11.73	33.9	ng/ul	6127760	2	10.25	3612190	20.0
n-Hexadecanoic acid	17.82	5.4	ng/ul	1373410	4	16.89	5049660	20.0
(DEL) Alkane: Cyc...	20.84	2.8	ng/ul	707229	5	21.10	5144570	20.0
(DEL) Alkane: Str...	22.14	2.0	ng/ul	520604	5	21.10	5144570	20.0
(DEL) Alkane: Str...	22.63	2.7	ng/ul	691386	6	23.24	5116440	20.0
(DEL) Alkane: Str...	23.16	3.0	ng/ul	775769	6	23.24	5116440	20.0
(DEL) Alkane: Str...	23.77	3.8	ng/ul	958729	6	23.24	5116440	20.0